

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112118\
 Data File : BF110936.D
 Acq On : 21 Nov 2018 18:52
 Operator : JU/SJ
 Sample : J5996-25 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AOC1-06-A

Manual Integrations
 APPROVED

mohammad
 11/23/2018 11:49:13 AM

Quant Time: Nov 23 02:00:32 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	62764	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	240209	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	97299	20.00	ng	-0.01
64) Phenanthrene-d10	11.48	188	161209	20.00	ng	0.00
76) Chrysene-d12	14.13	240	111876	20.00	ng	0.00
87) Perylene-d12	15.67	264	77116	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	125262	32.42	ng	0.00
7) Phenol-d6	6.56	99	247303	50.05	ng	-0.01
23) Nitrobenzene-d5	7.51	82	163818	39.27	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	10053	10.25	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	294604	43.24	ng	-0.01
79) Terphenyl-d14	13.06	244	252065	42.19	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	11949	2.086	ng	93
50) Dimethylphthalate	9.69	163	19011	2.619	ng	97
75) Fluoranthene	12.70	202	49944	5.768	ng	98
78) Pyrene	12.93	202	50929	5.912	ng	96
81) Benzo(a)anthracene	14.13	228	23769	3.555	ng	# 91
83) Chrysene	14.16	228	26917	4.225	ng	# 87
86) Indeno(1,2,3-cd)pyrene	17.28	276	11667	2.552	ng	# 91
88) Benzo(b)fluoranthene	15.22	252	27449m	5.950	ng	
90) Benzo(a)pyrene	15.61	252	17401	4.151	ng	# 69
92) Benzo(g,h,i)perylene	17.77	276	14624m	4.155	ng	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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